

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061318\
 Data File : BF106396.D
 Acq On : 13 Jun 2018 17:32
 Operator : JU/SJ
 Sample : J3246-01 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-02-061218

Manual Integrations
 APPROVED

Sohil
 6/14/2018 2:47:43 PM

Quant Time: Jun 14 03:27:52 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 13 10:51:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	159168	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	626825	20.00	ng	-0.01
38) Acenaphthene-d10	10.01	164	307173	20.00	ng	-0.01
63) Phenanthrene-d10	11.50	188	566862	20.00	ng	-0.01
75) Chrysene-d12	14.15	240	434903	20.00	ng	-0.01
86) Perylene-d12	15.69	264	426902	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	381650	40.58	ng	-0.01
7) Phenol-d6	6.61	99	505177	42.52	ng	-0.01
23) Nitrobenzene-d5	7.52	82	313572	29.43	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	127349	43.09	ng	-0.02
44) 2-Fluorobiphenyl	9.32	172	620395	25.84	ng	-0.01
78) Terphenyl-d14	13.08	244	590526	33.73	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.71	163	104895	4.780	ng	99
70) Phenanthrene	11.52	178	204917	7.381	ng	98
71) Anthracene	11.57	178	56664	2.037	ng	99
74) Fluoranthene	12.71	202	326939	11.043	ng	98
77) Pyrene	12.94	202	345160	12.677	ng	99
80) Benzo(a)anthracene	14.14	228	194672m	7.522	ng	
82) Chrysene	14.18	228	191224	8.093	ng	98
85) Indeno(1,2,3-cd)pyrene	17.25	276	98149	5.205	ng	# 90
87) Benzo(b)fluoranthene	15.24	252	273925m	10.615	ng	
88) Benzo(k)fluoranthene	15.26	252	91823m	3.593	ng	
89) Benzo(a)pyrene	15.62	252	195850	8.438	ng	99
91) Benzo(a,h,i)perylene	17.73	276	88601	4.711	ng	# 93

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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